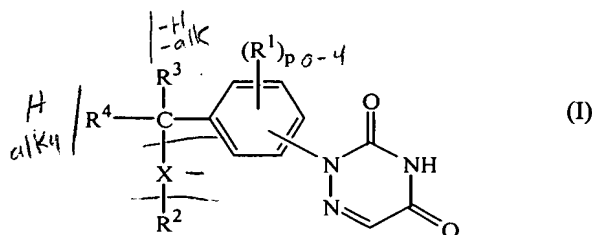


Note: Applicant uses:

- ~~Cross out text to indicate deletions~~
- Underline text to indicate additions

**Claims:**

1. (Currently Amended) A compound having the formula



a *N*-oxide, a pharmaceutically acceptable addition salt or a stereochemically isomeric form thereof, wherein:

*p* represents an integer being 0, 1, 2, 3 or 4;

*X* represents O, S, NR<sup>5</sup> or a direct bond;

*Y* represents O, S, NR<sup>5</sup>, or S(O)<sub>2</sub>;

each R<sup>1</sup> independently represents C<sub>1-6</sub>alkyl, halo, polyhaloC<sub>1-6</sub>alkyl, hydroxy, mercapto, C<sub>1-6</sub>alkyloxy, C<sub>1-6</sub>alkylthio, C<sub>1-6</sub>alkylcarbonyloxy, aryl, cyano, nitro, Het<sup>3</sup>, R<sup>6</sup>, NR<sup>7</sup>R<sup>8</sup> or C<sub>1-4</sub>alkyl substituted with Het<sup>3</sup>, R<sup>6</sup> or NR<sup>7</sup>R<sup>8</sup>;

R<sup>2</sup> represents Het<sup>1</sup>, C<sub>3-7</sub>cycloalkyl, C<sub>1-6</sub>alkyl or C<sub>1-6</sub>alkyl substituted with one or two substituents selected from hydroxy, cyano, amino, mono- or di(C<sub>1-4</sub>alkyl)amino, C<sub>1-6</sub>alkyloxy, C<sub>1-6</sub>alkylsulfonyloxy, C<sub>1-6</sub>alkyloxycarbonyl, C<sub>3-7</sub>cycloalkyl, aryl, aryloxy, arylthio, Het<sup>1</sup>, Het<sup>1</sup>oxy and Het<sup>1</sup>thio; and if *X* is O, S or NR<sup>5</sup>, then R<sup>2</sup> may also represent aminocarbonyl, aminothiocarbonyl, C<sub>1-4</sub>alkylcarbonyl, C<sub>1-4</sub>alkylthiocarbonyl, arylcarbonyl, arylthiocarbonyl, Het<sup>1</sup>carbonyl or Het<sup>1</sup>thiocarbonyl;

R<sup>3</sup> represents hydrogen, C<sub>1-6</sub>alkyl or C<sub>3-7</sub>cycloalkyl;

R<sup>4</sup> represents hydrogen, C<sub>1-6</sub>alkyl or C<sub>3-7</sub>cycloalkyl; or

R<sup>3</sup> and R<sup>4</sup> taken together form a C<sub>2-6</sub>alkanediyl;

R<sup>5</sup> represents hydrogen or C<sub>1-4</sub>alkyl;

D<sup>2</sup>  
each R<sup>6</sup> independently represents C<sub>1-6</sub>alkylsulfonyl, aminosulfonyl, mono- or di(C<sub>1-4</sub>alkyl)aminosulfonyl, mono- or di(benzyl)aminosulfonyl, polyhaloC<sub>1-6</sub>alkylsulfonyl, C<sub>1-6</sub>alkylsulfinyl, phenylC<sub>1-4</sub>alkylsulfonyl, piperazinylsulfonyl, aminopiperidinylsulfonyl, piperidinylaminosulfonyl, *N*-C<sub>1-4</sub>alkyl-*N*-piperidinylaminosulfonyl or mono- or di(C<sub>1-4</sub>alkyl)aminoC<sub>1-4</sub>alkylsulfonyl;

each R<sup>7</sup> and each R<sup>8</sup> are independently selected from the group consisting of: hydrogen, C<sub>1-4</sub>alkyl, hydroxyC<sub>1-4</sub>alkyl, dihydroxyC<sub>1-4</sub>alkyl, aryl, arylC<sub>1-4</sub>alkyl, C<sub>1-4</sub>alkyloxyC<sub>1-4</sub>alkyl, C<sub>1-4</sub>alkylcarbonyl, aminocarbonyl, arylcarbonyl, Het<sup>3</sup>carbonyl, C<sub>1-4</sub>alkylcarbonyloxy-C<sub>1-4</sub>alkylcarbonyl, hydroxyC<sub>1-4</sub>alkylcarbonyl, C<sub>1-4</sub>alkyloxy-carbonylcarbonyl, mono- or di(C<sub>1-4</sub>alkyl)aminoC<sub>1-4</sub>alkyl, arylaminocarbonyl, arylaminothiocarbonyl, Het<sup>3</sup>aminocarbonyl, Het<sup>3</sup>aminothiocarbonyl, C<sub>3-7</sub>cycloalkyl, pyridinylC<sub>1-4</sub>alkyl, C<sub>1-4</sub>alkanediyl-C(=O)-O-R<sup>14</sup>, -C(=O)-O-R<sup>14</sup>, -Y-C<sub>1-4</sub>alkanediyl-C(=O)-O-R<sup>14</sup>, Het<sup>3</sup>, Het<sup>4</sup> and R<sup>6</sup>;

R<sup>9</sup> and R<sup>10</sup> are each independently selected from the group consisting of: hydrogen, C<sub>1-4</sub>alkyl, hydroxyC<sub>1-4</sub>alkyl, dihydroxyC<sub>1-4</sub>alkyl, phenyl, phenylC<sub>1-4</sub>alkyl, C<sub>1-4</sub>alkyloxyC<sub>1-4</sub>alkyl, C<sub>1-4</sub>alkylcarbonyl, aminocarbonyl, phenylcarbonyl, Het<sup>3</sup>carbonyl, C<sub>1-4</sub>alkylcarbonyloxyC<sub>1-4</sub>alkylcarbonyl, hydroxyC<sub>1-4</sub>alkylcarbonyl, C<sub>1-4</sub>alkyloxy-carbonylcarbonyl, mono- or di(C<sub>1-4</sub>alkyl)aminoC<sub>1-4</sub>alkyl, phenylaminocarbonyl, phenylaminothiocarbonyl, Het<sup>3</sup>aminocarbonyl, Het<sup>3</sup>aminothiocarbonyl, C<sub>3-7</sub>cycloalkyl, pyridinylC<sub>1-4</sub>alkyl, C<sub>1-4</sub>alkanediyl-C(=O)-O-R<sup>14</sup>, -C(=O)-O-R<sup>14</sup>, -Y-C<sub>1-4</sub>alkanediyl-C(=O)-O-R<sup>14</sup>, Het<sup>3</sup>, Het<sup>4</sup> and R<sup>6</sup>;

each R<sup>11</sup> independently being selected from the group consisting of: hydroxy, mercapto, cyano, nitro, halo, trihalomethyl, C<sub>1-4</sub>alkyloxy, formyl, trihaloC<sub>1-4</sub>alkylsulfonyloxy, R<sup>6</sup>, NR<sup>7</sup>R<sup>8</sup>, C(=O)NR<sup>7</sup>R<sup>8</sup>, -C(=O)-O-R<sup>14</sup>, -Y-C<sub>1-4</sub>alkanediyl-C(=O)-O-R<sup>14</sup>, aryl, aryloxy, arylcarbonyl, C<sub>3-7</sub>cycloalkyl, C<sub>3-7</sub>cycloalkyloxy, phthalimide-2-yl, Het<sup>3</sup> and C(=O)Het<sup>3</sup>;

R<sup>12</sup> and R<sup>13</sup> are each independently selected from the group consisting of: hydrogen, C<sub>1-4</sub>alkyl, hydroxyC<sub>1-4</sub>alkyl, dihydroxyC<sub>1-4</sub>alkyl, phenyl, phenylC<sub>1-4</sub>alkyl, C<sub>1-4</sub>alkyloxyC<sub>1-4</sub>alkyl, C<sub>1-4</sub>alkylcarbonyl, phenylcarbonyl, C<sub>1-4</sub>alkylcarbonyloxyC<sub>1-4</sub>alkylcarbonyl, hydroxyC<sub>1-4</sub>alkylcarbonyl, C<sub>1-4</sub>alkyloxy-carbonylcarbonyl, mono- or di(C<sub>1-4</sub>alkyl)aminoC<sub>1-4</sub>alkyl, phenylaminocarbonyl, phenylaminothiocarbonyl, C<sub>3-</sub>

7cycloalkyl, pyridinylC<sub>1-4</sub>alkyl, C<sub>1-4</sub>alkanediyl-C(=O)-O-R<sup>14</sup>, -C(=O)-O-R<sup>14</sup>, -Y-C<sub>1-4</sub>alkanediyl-C(=O)-O-R<sup>14</sup> and R<sup>6</sup>;

each R<sup>14</sup> independently represents hydrogen, C<sub>1-4</sub>alkyl, C<sub>3-7</sub>cycloalkyl, aminocarbonylmethylene or mono-or di(C<sub>1-4</sub>alkyl)aminocarbonylmethylene;

aryl represents phenyl optionally substituted with one, two or three substituents each

independently selected from nitro, azido, cyano, halo, hydroxy, C<sub>1-4</sub>alkyl, C<sub>3-7</sub>cycloalkyl, C<sub>1-4</sub>alkyloxy, formyl, polyhaloC<sub>1-4</sub>alkyl, NR<sup>9</sup>R<sup>10</sup>, C(=O)NR<sup>9</sup>R<sup>10</sup>, C(=O)-O-R<sup>14</sup>, R<sup>6</sup>, -O-R<sup>6</sup>, phenyl, Het<sup>3</sup>, C(=O)Het<sup>3</sup> and C<sub>1-4</sub>alkyl substituted with hydroxy, C<sub>1-4</sub>alkyloxy, C(=O)-O-R<sup>14</sup>, -Y-C<sub>1-4</sub>alkanediyl-C(=O)-O-R<sup>14</sup>, Het<sup>3</sup> or NR<sup>9</sup>R<sup>10</sup>;

Het<sup>1</sup> represents a heterocycle selected from **the group consisting of:** pyrrolyl, pyrrolinyl, imidazolyl, imidazolynyl, pyrazolyl, pyrazolynyl, triazolyl, tetrazolyl, furanyl, tetrahydrofuranyl, thienyl, thiolanyl, dioxolanyl, oxazolyl, oxazolynyl, isoxazolyl, thiazolyl, thiazolynyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl, pyrrolidinyl, piperidinyl, piperazinyl, morpholinyl, thiomorpholinyl, dioxanyl, dithianyl, trithianyl, triazinyl, benzothienyl, isobenzothienyl, benzofuranyl, isobenzofuranyl, benzothiazolyl, benzoxazolyl, indolyl, isoindolyl, indolinyl, purinyl, 1H-pyrazolo[3,4-d]pyrimidinyl, benzimidazolyl, quinolyl, isoquinolyl, cinnolinyl, phtalazinyl, quinazolinyl, quinoxalinyl, thiazolopyridinyl, oxazolopyridinyl and imidazo[2,1-b]thiazolyl; wherein said heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from Het<sup>2</sup>, R<sup>11</sup> and C<sub>1-4</sub>alkyl optionally substituted with one or two substituents independently selected from Het<sup>2</sup> and R<sup>11</sup>;

**provided Het<sup>1</sup> is other than 2-substituted-pyridin-5-yl;**

Het<sup>2</sup> represents a heterocycle selected from **the group consisting of:** pyrrolyl, pyrrolinyl, imidazolyl, imidazolynyl, pyrazolyl, pyrazolynyl, triazolyl, tetrazolyl, furanyl, tetrahydrofuranyl, thienyl, thiolanyl, dioxolanyl, oxazolyl, oxazolynyl, isoxazolyl, thiazolyl, thiazolynyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl, dioxanyl, dithianyl, trithianyl, triazinyl, benzothienyl, isobenzothienyl, benzofuranyl, isobenzofuranyl, benzothiazolyl, benzoxazolyl, indolyl, isoindolyl, indolinyl, purinyl, 1H-pyrazolo[3,4-d]pyrimidinyl, benzimidazolyl, quinolyl,

D<sup>2</sup>

isoquinolyl, cinnolinyl, phtalazinyl, quinazoliny, quinoxaliny, thiazolopyridiny, oxazolopyridiny and imidazo[2,1-b]thiazoly; wherein said heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from Het<sup>4</sup>, R<sup>11</sup> and C<sub>1-4</sub>alkyl optionally substituted with one or two substituents independently selected from Het<sup>4</sup> and R<sup>11</sup>;

Het<sup>3</sup> represents a monocyclic heterocycle selected the group consisting of: from pyrrolidiny, piperidiny, piperaziny, morpholiny, thiomorpholiny and tetrahydropyrany; wherein said monocyclic heterocycles each independently may optionally be substituted with, where possible, one, two, three or four substituents each independently selected from hydroxy, C<sub>1-4</sub>alkyl, C<sub>1-4</sub>alkyloxy, C<sub>1-4</sub>alkylcarbonyl, piperidiny, NR<sup>12</sup>R<sup>13</sup>, C(=O)-O-R<sup>14</sup>, R<sup>6</sup> and C<sub>1-4</sub>alkyl substituted with one or two substituents independently selected from hydroxy, C<sub>1-4</sub>alkyloxy, phenyl, C(=O)-O-R<sup>14</sup>, -Y-C<sub>1-4</sub>alkanediyl-C(=O)-O-R<sup>14</sup>, R<sup>6</sup> and NR<sup>12</sup>R<sup>13</sup>;

Het<sup>4</sup> represents a monocyclic heterocycle selected from the group consisting of: pyrrolyl, imidazolyl, pyrazolyl, triazolyl, tetrazolyl, furanyl, thienyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridiny, pyrimidiny, pyraziny, pyranly, pyridaziny and triaziny.

2. (Currently Amended) A compound as claimed in claim 1 wherein:  
each R<sup>7</sup> and each R<sup>8</sup> are independently selected from the group consisting of: hydrogen, C<sub>1-4</sub>alkyl, hydroxyC<sub>1-4</sub>alkyl, dihydroxyC<sub>1-4</sub>alkyl, aryl, arylC<sub>1-4</sub>alkyl, C<sub>1-4</sub>alkyloxyC<sub>1-4</sub>alkyl, C<sub>1-4</sub>alkylcarbonyl, aminocarbonyl, arylcarbonyl, Het<sup>3</sup>carbonyl, C<sub>1-4</sub>alkylcarbonyloxy-C<sub>1-4</sub>alkylcarbonyl, hydroxyC<sub>1-4</sub>alkylcarbonyl, C<sub>1-4</sub>alkyloxy-carbonylcarbonyl, mono- or di(C<sub>1-4</sub>alkyl)aminoC<sub>1-4</sub>alkyl, arylaminocarbonyl, arylaminothiocarbonyl, Het<sup>3</sup>aminocarbonyl, Het<sup>3</sup>aminothiocarbonyl, C<sub>3-7</sub>cycloalkyl, pyridinyC<sub>1-4</sub>alkyl, C<sub>1-4</sub>alkanediyl-C(=O)-O-R<sup>14</sup>, -C(=O)-O-R<sup>14</sup>, -Y-C<sub>1-4</sub>alkanediyl-C(=O)-O-R<sup>14</sup>, Het<sup>3</sup> and R<sup>6</sup>;
- R<sup>9</sup> and R<sup>10</sup> are each independently selected from the group consisting of: hydrogen, C<sub>1-4</sub>alkyl, hydroxyC<sub>1-4</sub>alkyl, dihydroxyC<sub>1-4</sub>alkyl, phenyl, phenylC<sub>1-4</sub>alkyl, C<sub>1-4</sub>alkyloxyC<sub>1-4</sub>alkyl, C<sub>1-4</sub>alkylcarbonyl, aminocarbonyl, phenylcarbonyl,

12 Het<sup>3</sup>carbonyl, C<sub>1-4</sub>alkylcarbonyloxyC<sub>1-4</sub>alkylcarbonyl, hydroxyC<sub>1-4</sub>alkylcarbonyl, C<sub>1-4</sub>alkyloxycarbonylcarbonyl, mono- or di(C<sub>1-4</sub>alkyl)aminoC<sub>1-4</sub>alkyl, phenylaminocarbonyl, phenylaminothiocarbonyl, Het<sup>3</sup>aminocarbonyl, Het<sup>3</sup>aminothiocarbonyl, C<sub>3-7</sub>cycloalkyl, pyridinylC<sub>1-4</sub>alkyl, C<sub>1-4</sub>alkanediyl-C(=O)-O-R<sup>14</sup>, -C(=O)-O-R<sup>14</sup>, -Y-C<sub>1-4</sub>alkanediyl-C(=O)-O-R<sup>14</sup>, Het<sup>3</sup> and R<sup>6</sup>;

R<sup>11</sup> is being selected from the group consisting of: hydroxy, mercapto, cyano, nitro, halo, trihalomethyl, C<sub>1-4</sub>alkyloxy, formyl, trihaloC<sub>1-4</sub>alkylsulfonyloxy, R<sup>6</sup>, NR<sup>7</sup>R<sup>8</sup>, C(=O)NR<sup>7</sup>R<sup>8</sup>, -C(=O)-O-R<sup>14</sup>, -Y-C<sub>1-4</sub>alkanediyl-C(=O)-O-R<sup>14</sup>, aryl, aryloxy, arylcarbonyl, C<sub>3-7</sub>cycloalkyl, C<sub>3-7</sub>cycloalkyloxy, phthalimide-2-yl, Het<sup>3</sup>, Het<sup>4</sup> and C(=O)Het<sup>3</sup>; and

Het<sup>2</sup> represents a heterocycle selected from the group consisting of: pyrrolyl, pyrrolinyl, imidazolyl, imidazoliny, pyrazolyl, pyrazolinyl, triazolyl, tetrazolyl, furanyl, tetrahydrofuranyl, thienyl, thiolanyl, dioxolanyl, oxazolyl, oxazolinyl, isoxazolyl, thiazolyl, thiazolinyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl, dioxanyl, dithianyl, trithianyl, triazinyl, benzothienyl, isobenzothienyl, benzofuranyl, isobenzofuranyl, benzothiazolyl, benzoxazolyl, indolyl, isoindolyl, indolinyl, purinyl, 1H-pyrazolo[3,4-d]pyrimidinyl, benzimidazolyl, quinolyl, isoquinolyl, cinnolinyl, phtalazinyl, quinazolinyl, quinoxalinyl, thiazolopyridinyl, oxazolopyridinyl and imidazo[2,1-b]thiazolyl; wherein said heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from R<sup>11</sup> and C<sub>1-4</sub>alkyl optionally substituted with one or two substituents independently selected from R<sup>11</sup>.

3-9. (Canceled)

3

10. (Currently Amended) A compound of the formula: as claimed in claim 1 wherein the compound is

2-[3,5-dichloro-4-[1-methyl-1-(4-phenyl-2-thiazolyl)ethyl]phenyl]-1,2,4-triazine-3,5(2H,4H)-dione;

- 02
- 2-[3,5-dichloro-4-[1-[4-(3-chlorophenyl)-5-methyl-2-thiazolyl]-1-methylethyl]-phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione;  
2-[3,5-dichloro-4-[1-methyl-1-(5-phenyl-1,2,4-oxadiazol-3-yl)ethyl]phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione;  
2-[3,5-dichloro-4-[1-(4,5-diphenyl-2-thiazolyl)-1-methylethyl]phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione;  
2-[3,5-dichloro-4-[1-methyl-1-[5-(2-methylphenyl)-1,2,4-oxadiazol-3-yl]ethyl]-phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione;  
2-[3,5-dichloro-4-[1-methyl-1-(4-methyl-5-phenyl-2-thiazolyl)ethyl]phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione;  
2-[3,5-dichloro-4-[1-methyl-1-[4-phenyl-5-(3-pyridinyl)-2-thiazolyl]ethyl]phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione;  
2-[3,5-dichloro-4-[1-methyl-1-[4-phenyl-5-(phenylmethyl)-2-thiazolyl]ethyl]-phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione;  
2-[3,5-dichloro-4-[1-methyl-1-[5-(4-pyridinyl)-1,2,4-oxadiazol-3-yl]ethyl]phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione;  
2-[3,5-dichloro-4-[1-methyl-1-[4-(3-thienyl)-2-thiazolyl]ethyl]phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione;  
2-[3,5-dichloro-4-[1-[4-(2-furanyl)-2-thiazolyl]-1-methylethyl]phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione;  
2-[3,5-dichloro-4-[1-methyl-1-[5-(3-pyridinyl)-1,2,4-oxadiazol-3-yl]ethyl]phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione;  
2-[3,5-dichloro-4-[1-methyl-1-[5-(2-methyl-3-pyridinyl)-1,2,4-oxadiazol-3-yl]ethyl]phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione; or  
2-[3,5-dichloro-4-[1-methyl-1-(5-phenyl-1,3,4-oxadiazol-2-yl)ethyl]phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione; or a *N*-oxide, a pharmaceutically acceptable addition salt or a stereochemically isomeric form thereof.

4  
11. (Previously Amended) A composition comprising a pharmaceutically acceptable carrier and, as active ingredient, a therapeutically effective amount of a compound as claimed in claim 1.

12. through 18. (Previously Canceled)

19. through 55. (Currently Canceled)

5  
56.

02 (New) A compound as claimed in claim 1 provided that in those compounds wherein X is a direct bond, at least one of R<sup>3</sup> and R<sup>4</sup> is hydrogen, and R<sup>2</sup> is 3-pyridinyl optionally substituted in the 6 position with an optionally substituted alkyl or acyl group are excluded.

6  
57.

(New) A compound as claimed in claim 2 provided that in those compounds wherein X is a direct bond, at least one of R<sup>3</sup> and R<sup>4</sup> is hydrogen, and R<sup>2</sup> is 3-pyridinyl optionally substituted in the 6 position with an optionally substituted alkyl or acyl group are excluded.

7  
58.

(New) A compound as claimed in claim 1 wherein the 6-azauracil moiety is in the para position relative to the carbon atom bearing the -X-R<sup>2</sup>, R<sup>3</sup> and R<sup>4</sup> substituents.

8  
59.

(New) A compound as claimed in claim 2 wherein the 6-azauracil moiety is in the para position relative to the carbon atom bearing the -X-R<sup>2</sup>, R<sup>3</sup> and R<sup>4</sup> substituents.

9  
60.

5 (New) A compound as claimed in claim 56 wherein the 6-azauracil moiety is in the para position relative to the carbon atom bearing the -X-R<sup>2</sup>, R<sup>3</sup> and R<sup>4</sup> substituents.

10  
61.

16 (New) A compound as claimed in claim 57 wherein the 6-azauracil moiety is in the para position relative to the carbon atom bearing the -X-R<sup>2</sup>, R<sup>3</sup> and R<sup>4</sup> substituents.

11  
62.

(New) A compound as claimed in claim 1 wherein R<sup>2</sup> is a monocyclic heterocycle selected from the group consisting of: pyrrolyl, imidazolyl, pyrazolyl, triazolyl, tetrazolyl, furanyl, thienyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranlyl, pyridazinyl and triazinyl, wherein said monocyclic heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from Het<sup>2</sup>, R<sup>11</sup> and C<sub>1-4</sub>alkyl optionally substituted with Het<sup>2</sup> or R<sup>11</sup>.

12  
63.

12  
A compound as claimed in claim 2 wherein  $R^2$  is a monocyclic heterocycle selected from the group consisting of: pyrrolyl, imidazolyl, pyrazolyl, triazolyl, tetrazolyl, furanyl, thienyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl and triazinyl, wherein said monocyclic heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from  $Het^2$ ,  $R^{11}$  and  $C_{1-4}alkyl$  optionally substituted with  $Het^2$  or  $R^{11}$ .

13  
64.

13  
A compound as claimed in claim 58 wherein  $R^2$  is a monocyclic heterocycle selected from the group consisting of: pyrrolyl, imidazolyl, pyrazolyl, triazolyl, tetrazolyl, furanyl, thienyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl and triazinyl, wherein said monocyclic heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from  $Het^2$ ,  $R^{11}$  and  $C_{1-4}alkyl$  optionally substituted with  $Het^2$  or  $R^{11}$ .

14  
65.

14  
A compound as claimed in claim 59 wherein  $R^2$  is a monocyclic heterocycle selected from the group consisting of: pyrrolyl, imidazolyl, pyrazolyl, triazolyl, tetrazolyl, furanyl, thienyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl and triazinyl, wherein said monocyclic heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from  $Het^2$ ,  $R^{11}$  and  $C_{1-4}alkyl$  optionally substituted with  $Het^2$  or  $R^{11}$ .

15  
66.

15  
A compound as claimed in claim 60 wherein  $R^2$  is a monocyclic heterocycle selected from the group consisting of: pyrrolyl, imidazolyl, pyrazolyl, triazolyl, tetrazolyl, furanyl, thienyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl and triazinyl, wherein said monocyclic heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from  $Het^2$ ,  $R^{11}$  and  $C_{1-4}alkyl$  optionally substituted with  $Het^2$  or  $R^{11}$ .



<sup>16</sup>  
~~67.~~ <sup>8</sup>  
D<sup>2</sup> A compound as claimed in claim ~~59~~ wherein R<sup>2</sup> is a monocyclic heterocycle selected from the group consisting of: pyrrolyl, imidazolyl, pyrazolyl, triazolyl, tetrazolyl, furanyl, thienyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl and triazinyl, wherein said monocyclic heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from Het<sup>2</sup>, R<sup>11</sup> and C<sub>1-4</sub>alkyl optionally substituted with Het<sup>2</sup> or R<sup>11</sup>.

<sup>17</sup>  
~~68.~~ <sup>9</sup>  
A compound as claimed in claim ~~60~~ wherein R<sup>2</sup> is a monocyclic heterocycle selected from the group consisting of: pyrrolyl, imidazolyl, pyrazolyl, triazolyl, tetrazolyl, furanyl, thienyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl and triazinyl, wherein said monocyclic heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from Het<sup>2</sup>, R<sup>11</sup> and C<sub>1-4</sub>alkyl optionally substituted with Het<sup>2</sup> or R<sup>11</sup>.

<sup>18</sup>  
~~69.~~ <sup>10</sup>  
A compound as claimed in claim ~~61~~ wherein R<sup>2</sup> is a monocyclic heterocycle selected from the group consisting of: pyrrolyl, imidazolyl, pyrazolyl, triazolyl, tetrazolyl, furanyl, thienyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl and triazinyl, wherein said monocyclic heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from Het<sup>2</sup>, R<sup>11</sup> and C<sub>1-4</sub>alkyl optionally substituted with Het<sup>2</sup> or R<sup>11</sup>.

<sup>19</sup>  
~~70.~~ (New) A compound as claimed in claim ~~1~~ wherein R<sup>3</sup> and R<sup>4</sup> are both methyl and -X-R<sup>2</sup> is Het<sup>1</sup>.

<sup>20</sup>  
~~71.~~ (New) A compound as claimed in claim ~~2~~ wherein R<sup>3</sup> and R<sup>4</sup> are both methyl and -X-R<sup>2</sup> is Het<sup>1</sup>.

<sup>21</sup>  
~~72.~~ (New) A compound as claimed in claim ~~56~~<sup>5</sup> wherein R<sup>3</sup> and R<sup>4</sup> are both methyl and -X-R<sup>2</sup> is Het<sup>1</sup>.

<sup>22</sup>  
~~73.~~ (New) A compound as claimed in claim ~~57~~<sup>6</sup> wherein R<sup>3</sup> and R<sup>4</sup> are both methyl and -X-R<sup>2</sup> is Het<sup>1</sup>.

<sup>23</sup>  
~~74.~~ (New) A compound as claimed in claim ~~58~~<sup>7</sup> wherein R<sup>3</sup> and R<sup>4</sup> are both methyl and -X-R<sup>2</sup> is Het<sup>1</sup>.

<sup>24</sup>  
~~75.~~ (New) A compound as claimed in claim ~~59~~<sup>8</sup> wherein R<sup>3</sup> and R<sup>4</sup> are both methyl and -X-R<sup>2</sup> is Het<sup>1</sup>.

<sup>25</sup>  
~~76.~~ (New) A compound as claimed in claim ~~60~~<sup>9</sup> wherein R<sup>3</sup> and R<sup>4</sup> are both methyl and -X-R<sup>2</sup> is Het<sup>1</sup>.

<sup>26</sup>  
~~77.~~ (New) A compound as claimed in claim ~~61~~<sup>10</sup> wherein R<sup>3</sup> and R<sup>4</sup> are both methyl and -X-R<sup>2</sup> is Het<sup>1</sup>.

<sup>27</sup>  
~~78.~~ (New) A compound as claimed in claim ~~62~~<sup>11</sup> wherein R<sup>3</sup> and R<sup>4</sup> are both methyl and -X-R<sup>2</sup> is Het<sup>1</sup>.

<sup>28</sup>  
~~79.~~ (New) A compound as claimed in claim ~~63~~<sup>12</sup> wherein R<sup>3</sup> and R<sup>4</sup> are both methyl and -X-R<sup>2</sup> is Het<sup>1</sup>.

<sup>29</sup>  
~~80.~~ (New) A compound as claimed in claim ~~64~~<sup>13</sup> wherein R<sup>3</sup> and R<sup>4</sup> are both methyl and -X-R<sup>2</sup> is Het<sup>1</sup>.

<sup>30</sup>  
~~81.~~ (New) A compound as claimed in claim ~~65~~<sup>14</sup> wherein R<sup>3</sup> and R<sup>4</sup> are both methyl and -X-R<sup>2</sup> is Het<sup>1</sup>.

<sup>31</sup>  
~~82.~~ (New) A compound as claimed in claim ~~66~~<sup>15</sup> wherein R<sup>3</sup> and R<sup>4</sup> are both methyl and -X-R<sup>2</sup> is Het<sup>1</sup>.

<sup>32</sup>  
~~83.~~ (New) A compound as claimed in claim ~~67~~<sup>16</sup> wherein R<sup>3</sup> and R<sup>4</sup> are both methyl and -X-R<sup>2</sup> is Het<sup>1</sup>.

33  
84. (New) A compound as claimed in claim ~~88~~<sup>17</sup> wherein  $R^3$  and  $R^4$  are both methyl and  $-X-R^2$  is Het<sup>1</sup>.

34  
85. (New) A compound as claimed in claim ~~89~~<sup>18</sup> wherein  $R^3$  and  $R^4$  are both methyl and  $-X-R^2$  is Het<sup>1</sup>.

35  
86. (New) A compound as claimed in claim ~~90~~<sup>19</sup> wherein p is 1 or 2 and each  $R^1$  is chloro.

36  
87. (New) A compound as claimed in claim ~~91~~<sup>20</sup> wherein p is 1 or 2 and each  $R^1$  is chloro.

37  
88. (New) A compound as claimed in claim ~~92~~<sup>21</sup> wherein p is 1 or 2 and each  $R^1$  is chloro.

38  
89. (New) A compound as claimed in claim ~~93~~<sup>22</sup> wherein p is 1 or 2 and each  $R^1$  is chloro.

39  
90. (New) A compound as claimed in claim ~~94~~<sup>23</sup> wherein p is 1 or 2 and each  $R^1$  is chloro.

40  
91. (New) A compound as claimed in claim ~~95~~<sup>24</sup> wherein p is 1 or 2 and each  $R^1$  is chloro.

41  
92. (New) A compound as claimed in claim ~~96~~<sup>25</sup> wherein p is 1 or 2 and each  $R^1$  is chloro.

42  
93. (New) A compound as claimed in claim ~~97~~<sup>26</sup> wherein p is 1 or 2 and each  $R^1$  is chloro.

43  
94. (New) A compound as claimed in claim ~~98~~<sup>27</sup> wherein p is 1 or 2 and each  $R^1$  is chloro.

<sup>44</sup>  
~~95.~~ (New) A compound as claimed in claim ~~62~~<sup>12</sup> wherein p is 1 or 2 and each R<sup>1</sup> is chloro.

<sup>45</sup>  
~~96.~~ (New) A compound as claimed in claim ~~63~~<sup>13</sup> wherein p is 1 or 2 and each R<sup>1</sup> is chloro.

<sup>46</sup>  
~~97.~~ (New) A compound as claimed in claim ~~64~~<sup>14</sup> wherein p is 1 or 2 and each R<sup>1</sup> is chloro.

<sup>47</sup>  
~~98.~~ (New) A compound as claimed in claim ~~65~~<sup>15</sup> wherein p is 1 or 2 and each R<sup>1</sup> is chloro.

<sup>48</sup>  
~~99.~~ (New) A compound as claimed in claim ~~66~~<sup>16</sup> wherein p is 1 or 2 and each R<sup>1</sup> is chloro.

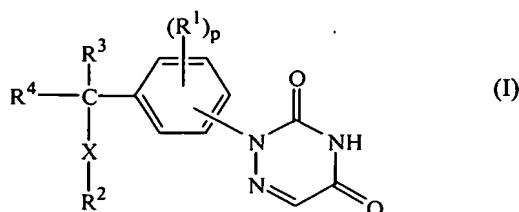
<sup>49</sup>  
~~100.~~ (New) A compound as claimed in claim ~~67~~<sup>17</sup> wherein p is 1 or 2 and each R<sup>1</sup> is chloro.

<sup>50</sup>  
~~101.~~ (New) A compound as claimed in claim ~~68~~<sup>18</sup> wherein p is 1 or 2 and each R<sup>1</sup> is chloro.

<sup>51</sup>  
~~102.~~ (New) A compound as claimed in claim ~~1~~ wherein R<sup>3</sup> and R<sup>4</sup> are both methyl, -X-R<sup>2</sup> is optionally substituted 2-thiazolyl or 3-oxadiazolyl, the 6-azauracil moiety is in the para position relative to the carbon atom bearing the -X-R<sup>2</sup>, R<sup>3</sup> and R<sup>4</sup> substituents, and p is 2 whereby both R<sup>1</sup> substituents are chloro positioned ortho relative to the carbon atom bearing the -X-R<sup>2</sup>, R<sup>3</sup> and R<sup>4</sup> substituents.

<sup>52</sup>  
~~103.~~ (New) A compound as claimed in claim ~~2~~ wherein R<sup>3</sup> and R<sup>4</sup> are both methyl, -X-R<sup>2</sup> is optionally substituted 2-thiazolyl or 3-oxadiazolyl, the 6-azauracil moiety is in the para position relative to the carbon atom bearing the -X-R<sup>2</sup>, R<sup>3</sup> and R<sup>4</sup> substituents, and p is 2 whereby both R<sup>1</sup> substituents are chloro positioned ortho relative to the carbon atom bearing the -X-R<sup>2</sup>, R<sup>3</sup> and R<sup>4</sup> substituents.

<sup>53</sup>  
~~104.~~ (New) A compound having the formula



a *N*-oxide, a pharmaceutically acceptable addition salt or a stereochemically isomeric form thereof, wherein :

*p* represents an integer being 0, 1 or 2;

*X* represents O, S, NR<sup>5</sup> or a direct bond;

*Y* represents O, S, NR<sup>5</sup>, or S(O)<sub>2</sub>;

each R<sup>1</sup> independently represents chloro or trifluoromethyl;

R<sup>2</sup> represents Het<sup>1</sup> or C<sub>1-6</sub>alkyl substituted with one or two substituents selected from

hydroxy, cyano, amino, mono- or di(C<sub>1-4</sub>alkyl)amino, C<sub>1-6</sub>alkyloxy,

C<sub>1-6</sub>alkylsulfonyloxy, C<sub>1-6</sub>alkyloxycarbonyl, C<sub>3-7</sub>cycloalkyl, aryl, aryloxy, arylthio, Het<sup>1</sup>,

Het<sup>1</sup>oxy and Het<sup>1</sup>thio; and if *X* is O, S or NR<sup>5</sup>, then R<sup>2</sup> may also represent

aminocarbonyl, aminothiocarbonyl, C<sub>1-4</sub>alkylcarbonyl, C<sub>1-4</sub>alkylthiocarbonyl,

arylcarbonyl, arylthiocarbonyl, Het<sup>1</sup>carbonyl or Het<sup>1</sup>thiocarbonyl;

R<sup>3</sup> represents hydrogen, C<sub>1-6</sub>alkyl or C<sub>3-7</sub>cycloalkyl;

R<sup>4</sup> represents hydrogen, C<sub>1-6</sub>alkyl or C<sub>3-7</sub>cycloalkyl; or R<sup>3</sup> and R<sup>4</sup> taken together form a C<sub>2-6</sub>alkanediyl;

R<sup>5</sup> represents hydrogen or C<sub>1-4</sub>alkyl;

each R<sup>6</sup> independently represents C<sub>1-6</sub>alkylsulfonyl, aminosulfonyl, mono- or

di(C<sub>1-4</sub>alkyl)aminosulfonyl, mono- or di(benzyl)aminosulfonyl,

polyhaloC<sub>1-6</sub>alkylsulfonyl, C<sub>1-6</sub>alkylsulfinyl, phenylC<sub>1-4</sub>alkylsulfonyl,

piperazinylsulfonyl, aminopiperidinylsulfonyl, piperidinylaminosulfonyl, *N*-C<sub>1-4</sub>alkyl-*N*-piperidinylaminosulfonyl or mono-or di(C<sub>1-4</sub>alkyl)aminoC<sub>1-4</sub>alkylsulfonyl;

each R<sup>7</sup> and each R<sup>8</sup> are independently selected from hydrogen, C<sub>1-4</sub>alkyl, hydroxyC<sub>1-4</sub>alkyl,

dihydroxyC<sub>1-4</sub>alkyl, aryl, arylC<sub>1-4</sub>alkyl, C<sub>1-4</sub>alkyloxyC<sub>1-4</sub>alkyl, C<sub>1-4</sub>alkylcarbonyl,

aminocarbonyl, arylcarbonyl, Het<sup>3</sup>carbonyl, C<sub>1-4</sub>alkylcarbonyloxy-C<sub>1-4</sub>alkylcarbonyl,

hydroxyC<sub>1-4</sub>alkylcarbonyl, C<sub>1-4</sub>alkyloxycarbonylcarbonyl, mono- or

di(C<sub>1-4</sub>alkyl)aminoC<sub>1-4</sub>alkyl, arylaminocarbonyl, arylaminothiocarbonyl,

Het<sup>3</sup>aminocarbonyl, Het<sup>3</sup>aminothiocarbonyl, C<sub>3-7</sub>cycloalkyl, pyridinylC<sub>1-4</sub>alkyl,

$C_{1-4}$ alkanediyl- $C(=O)-O-R^{14}$ ,  $-C(=O)-O-R^{14}$ ,  $-Y-C_{1-4}$ alkanediyl- $C(=O)-O-R^{14}$ ,  $Het^3$ ,  $Het^4$  and  $R^6$ ;

02  $R^9$  and  $R^{10}$  are each independently selected from hydrogen,  $C_{1-4}$ alkyl, hydroxy $C_{1-4}$ alkyl, dihydroxy $C_{1-4}$ alkyl, phenyl, phenyl $C_{1-4}$ alkyl,  $C_{1-4}$ alkyloxy $C_{1-4}$ alkyl,  $C_{1-4}$ alkylcarbonyl, aminocarbonyl, phenylcarbonyl,  $Het^3$ carbonyl,  $C_{1-4}$ alkylcarbonyloxy $C_{1-4}$ alkylcarbonyl, hydroxy $C_{1-4}$ alkylcarbonyl,  $C_{1-4}$ alkyloxy $C_{1-4}$ alkylcarbonyl, mono- or di( $C_{1-4}$ alkyl)amino $C_{1-4}$ alkyl, phenylaminocarbonyl, phenylaminothiocarbonyl,  $Het^3$ aminocarbonyl,  $Het^3$ aminothiocarbonyl,  $C_{3-7}$ cycloalkyl, pyridinyl $C_{1-4}$ alkyl,  $C_{1-4}$ alkanediyl- $C(=O)-O-R^{14}$ ,  $-C(=O)-O-R^{14}$ ,  $-Y-C_{1-4}$ alkanediyl- $C(=O)-O-R^{14}$ ,  $Het^3$ ,  $Het^4$  and  $R^6$ ;

each  $R^{11}$  independently being selected from hydroxy, cyano, nitro, halo,  $C_{1-4}$ alkyloxy, formyl,  $NR^7R^8$ ,  $C(=O)NR^7R^8$ ,  $-C(=O)-O-R^{14}$ , aryl, arylcarbonyl,  $Het^3$  and  $C(=O)Het^3$ ;

each  $R^{14}$  independently represents hydrogen,  $C_{1-4}$ alkyl,  $C_{3-7}$ cycloalkyl,

aminocarbonylmethylene or mono- or di( $C_{1-4}$ alkyl)aminocarbonylmethylene;

aryl represents phenyl optionally substituted with one, two or three substituents each independently selected from nitro, azido, cyano, halo, hydroxy,  $C_{1-4}$ alkyl,

$C_{3-7}$ cycloalkyl,  $C_{1-4}$ alkyloxy, formyl, polyhalo $C_{1-4}$ alkyl,  $NR^9R^{10}$ ,  $C(=O)NR^9R^{10}$ ,

$C(=O)-O-R^{14}$ ,  $R^6$ ,  $-O-R^6$ , phenyl,  $Het^3$ ,  $C(=O)Het^3$  and  $C_{1-4}$ alkyl substituted with hydroxy,  $C_{1-4}$ alkyloxy,  $C(=O)-O-R^{14}$ ,  $-Y-C_{1-4}$ alkanediyl- $C(=O)-O-R^{14}$ ,  $Het^3$  or  $NR^9R^{10}$ ;

$Het^1$  represents a heterocycle selected from pyrrolyl, imidazolyl, pyrazolyl, triazolyl, tetrazolyl, furanyl, thienyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl and triazinyl, wherein said monocyclic heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from  $Het^2$ ,  $R^{11}$  and  $C_{1-4}$ alkyl optionally substituted with  $Het^2$  or  $R^{11}$ ; provided  $Het^1$  is other than 2-substituted-pyridin-5-yl;

$Het^2$  represents a heterocycle selected from furanyl, thienyl, pyridinyl or benzothienyl,

wherein said aromatic heterocycles each independently may optionally be substituted

with one, or where possible, two or three substituents each independently selected from  $Het^4$ ,  $R^{11}$  and  $C_{1-4}$ alkyl optionally substituted with  $R^{11}$ ;

$Het^3$  represents a monocyclic heterocycle selected from piperidinyl, piperazinyl, morpholinyl and tetrahydropyranyl each independently and optionally substituted with, where possible, one, two, three or four substituents each independently selected from hydroxy,  $C_{1-4}$ alkyl,

D

C(=O)-O-R<sup>14</sup>, C<sub>1-4</sub>alkylcarbonyl, R<sup>6</sup>, piperidinyl and C<sub>1-4</sub>alkyl substituted with one or two substituents independently selected from hydroxy, C<sub>1-4</sub>alkyloxy, C(=O)-O-R<sup>14</sup> and phenyl; Het<sup>4</sup> represents a monocyclic heterocycle selected from thienyl or pyridinyl.

105. (New) A compound as claimed in claim 104, wherein when X is a direct bond, at least one of R<sup>3</sup> and R<sup>4</sup> is hydrogen, and R<sup>2</sup> is 3-pyridinyl, then R<sup>2</sup> is not substituted in the 6 position with an optionally substituted alkyl or acyl group.

106. (New) A compound as claimed in claim 104 wherein the 6-azauracil moiety is in the para position relative to the carbon atom bearing the -X-R<sup>2</sup>, R<sup>3</sup> and R<sup>4</sup> substituents.

107. (New) A compound as claimed in claim 104 wherein R<sup>2</sup> is a monocyclic heterocycle selected from the group consisting of: pyrrolyl, imidazolyl, pyrazolyl, triazolyl, tetrazolyl, furanyl, thienyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl and triazinyl, wherein said monocyclic heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from Het<sup>2</sup>, R<sup>11</sup> and C<sub>1-4</sub>alkyl optionally substituted with Het<sup>2</sup> or R<sup>11</sup>.

108. (New) A compound as claimed in claim 104 wherein R<sup>3</sup> and R<sup>4</sup> are both methyl and -X-R<sup>2</sup> is Het<sup>1</sup>.

109. (New) A compound as claimed in claim 104 wherein p is 1 or 2 and each R<sup>1</sup> is chloro.

110. (New) A compound as claimed in claim 104 wherein R<sup>3</sup> and R<sup>4</sup> are both methyl, -X-R<sup>2</sup> is optionally substituted 2-thiazolyl or 3-oxadiazolyl, the 6-azauracil moiety is in the para position relative to the carbon atom bearing the -X-R<sup>2</sup>, R<sup>3</sup> and R<sup>4</sup> substituents, and p is 2 whereby both R<sup>1</sup> substituents are chloro positioned ortho relative to the carbon atom bearing the -X-R<sup>2</sup>, R<sup>3</sup> and R<sup>4</sup> substituents.